S/058/61/000/010/069/100 A001/A101

24.7700

AUTHOR:

Perlin, Yu.Ye.

TITLE

On the theory of quantum transitions of a localized electron into polaron states

PERIODICAL: Referativnyy zhurnal. Fizika, no. 10, 1961, 237, abstract 10E22 ("Uch. zap. Kishinevsk. un-t", 1960, v. 55, 95 - 105)

TEXT: The author considers the process of thermoionization of F-centers in alkalihalide crystals accompanied by quantum transitions of electrons from the ground state of the F-center into polaron state. Non-adiabaticity operator is assumed as a perturbation causing the transition. An essential difference of the effective field in which polaron moves from the Coulomb field is taken into associat. It is assumed that kinetic energy of polaron is considerably less than the energy of phon, due to which circumstance quasi-classical wave functions of polaron are used in continuous spectrum. Quantum transitions are assumed only in the s-state of this spectrum. Probability of thermoionization W is calculated on assumption that electrons interact only with longitudinal optical oscillations, dispersion of oscillation frequencies being neglected. An explicit analytical

Card 1/2

S/058/61/000/010/069/100
A001/8/1

expression of W as a function of temperature has been found. In particular, at right compenstures W depends exponentially on inverse temperature. The forext tomential factor for KC1 and NaC1 crystals is of the order of 10 sec-1.

M. Krivoglaz

Abstractar's note; Complete translation]

S/058/61/000/012/025 0.3: A058/A101

AUTHOR:

Ferlin, Yu. Ye.

TITLE:

Concerning the probability of nonradiative transition in the discrete spectrum of impurity centers

PERIODICAL: Referativnyy zhurnal, Fizika, no. 12, 1961, 196, abstract 12V25. ("Uch.zap. Kishinevsk. un-t", 1960, no. 55, 107-112)

The probability of multiphonon nonradiative transition in an F-center TEXT: was calculated on the assumption that electrons interact strongly with the optical but weakly with the acoustic branches of lattice vibrations. Acoustic interaction is taken into account as a transition-generating perturbation. The obtained formulae enable one to numerically evaluate the transition probability if the electron wave functions of the impurity center are known and the temperature dependence of thermal transition has been analyzed.

[Abstracter's note: Complete translation]

Card 1/1

PERLIN, Yu. Ye.

Consideration of the polaron effect in the theory of multiphonon thermal ionization. Fiz. tver. tela 2 no.2: 242-254 F '60. (MIRA 14:8)

1. Kishimevskiy gosudarstvennyy universitet.
(Iomic crystals--Defects) (Thermionic emission)

CIA-RDP86-00513R001240110006-6 "APPROVED FOR RELEASE: 06/15/2000 84815 5/181/60/002/008/050/052/XX BOO6/BO70 TEXT: The aim of the author was to generalize the theory localized near solution to the case of electrons which are author assumes fluorescence of atoms to the with the phonon field the lattice defects and interact with the phonon field the case of a toms to the case of the phonon field the lattice defects and interact with the phonon field the lattice defects and interact with the phonon field the lattice defects and interact with the phonon field the lattice defects and interact with the phonon field the lattice defects and interact with the phonon field the lattice defects and interact with the phonon field the lattice defects and interact with the phonon field the lattice defects and interact with the phonon field the lattice defects and interact with the phonon field the lattice defects and interact with the phonon field the lattice defects and interact with the phonon field the lattice defects and interact with the phonon field the lattice defects and interact with the phonon field the lattice defects and interact with the phonon field the lattice defects and interact with the phonon field the lattice defects and interact with the phonon field the lattice defects and interact with the phonon field the lattice defects and interact with the phonon field the lattice defects are considered to the lattice defects and lattice defects are considered to the lattice defects are co Impurity Fluorescence in Crystals 31 fluorescence of atoms to the case of electrons which are author assumes phonon field mhe author are localized near phonon field a dielectric are lattice defects and interact with impurity center of a dielectric that the optical electrons of an impurity center of a dielectric are lattice defects and interact with impurity center of a dielectric are lattice defects and interact with impurity center of a dielectric are lattice defects and interact with impurity center of a dielectric are lattice defects. lattice defects and interact with the phonon of the substance The the optical electrons of the substance that the optical bound than the intrinsic electrons of the substance less attrongly 24.3500 that the optical electrons of the intrinsic impurities lies in a region of the less the light absorption band of the fore, less strongly bound than the intrinsic electrons of the substance of the method of the fore, the light absorption band of the pure crystal Using the method of spectrum which is transparent AUTHOR: fore, the light absorption band of the impurities lies in a region of the impurities lies in the method of the impurities lies form a "fast" absorption band of the pure crystal electrons as "function the intrinsic obtained as and the impurity electrons and the crystal is obtained and the spectrum which the crystal is impurity electrons and the adiabatic approximation, in to of the impurity electrons are adiabatic approximation variables of the impurity of the configuration variables of the impurity electrons and the subsystem, configuration variables of the impurity electrons and the subsystem, configuration variables of the impurity electrons and the subsystem, configuration variables of the configuration variables of the configuration the configuration the configuration variables of the configuration that the configuration the configuration the TITLE: PERIODICAL: subsystem, the Bamiltonian wariables The energy of the configuration with them of the energy of the configuration with them subsystem is electron of a lift the electron phonon of the configuration fixed configuration fixed configuration of the latter of the phonon dielectric for a energy of the latter lift the electron phonon dielectric for a energy of the latter latter of the potential energy of the latter latt

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Impurity Fluorescence in Crystals

Card 2/3

S/181/60/002/008/050/052/XX B006/B070

interaction is not small, the Hamiltonian \hat{H}^{O} can be treated again in adiabatic approximation in which the localized electrons can be considered to be a "fast" subsystem with respect to the lattice vibrations. The formulas for the approximate wave function of \hat{H}^{O} and the approximate eigenvalue of \hat{H}^0 were derived by S. ... Pekar in Ref. 1. The possibility of radiationless transitions between the states considered is ignored If the crystal is irradiated with light of a frequency inside the impurity absorption band, luminescence may appear. The appearance of this luminescence is described in three stages. The first and the third were considered by Pekar. This luminescence is regarded as phosphorescence. It can also be regarded as fluorescence and looked upon as a single quantum transition: A photon $\hbar\Omega_{c}$ is absorbed and, simultaneously, another photon $k\Omega$ is emitted. The excited states of electrons and lattice play here the role of virtual states. Now, generalizing Weisskopf's theory the author considers three types of states: the initial state \$\psi_1\$ of the impurity electrons, the wirtual state in which the impurity ions are in the excited state ψ_2 , and the final state in which the electrons are again in the state 12. The theoretical study of the fluorescence probability and the spectral characteristics of the fluorescence lead to the following con-

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Impurity Pluorescence in Crystals

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clusions: 1) The spectrum of fluorescence excitation coincides with that of the absorption of light; 2) A Stokes shift appears: The fluorescence maximum shows a red shift at $\Delta\Omega$ = ω a compared to the absorption maximum; 3) The fluorescence curve represents a reflection of the impurity absorption band in the straight line $\Omega = \Omega$. The theoretical results are compared with the experimental results (Ref. 9) on the infrared fluorescence of the F-centers in KCl (Table) and NaCl at 77°K. The author thanks Professor S. I. Pekar for interest and discussions, and Professor A. I. Ansel'n and Professor L. E. Gurevich for discussions L. N. Ovander is mentioned. There are 1 table and 9 references: 6 Soviet, 1 US, 1 German and 1 British.

ASSOCIATION: Kishinevskiy gosudarstvennyy universitet (Kishinev State University)

SUBMITTED: January 19, 1960

Card 3/3

S/181/60/002/008/051/052/XX B006/B070

2 4.3500 AUTHOR:

Perlin, Yu. Ye.

TITLE

Consideration of Non radiative Transitions in the Theory

of Impurity Fluorescence

PERIODICAL:

Fizika tverdogo tela, 1960. Vol 2, No 8, pp 1928-1935

TEXT: This paper is in continuation of two previous papers (Refs. 1.2) in which a study was made of the fluorescence in a crystal caused by its irradiation with light of a frequency in the impurity absorption band. It was assumed that the impurity absorption band lies in a spectral region which is transparent to the pure substance, and that the optical electrons of an impurity center interact strongly with the crystal lattice vibrations. The fluorescence was regarded as a single quantum transition (absorption of a photon Ω_0 and a simultaneous emission of a photon Ω_0 The mathematical method developed in Ref. 2 is a generalization of Weisskopf's perturbation theory to strong electron-photon interaction, and permits a calculation of the spectral distribution curve. However, the non-radiative transitions between the virtual excited states and the

Card 1/3

84816

Consideration of Non-radiative Transitions S/18:/60/002/008/05:/052/XX in the Theory of Impurity Fluorescence B006/B070

final state of the system were not taken into account and, therefore, the quantum yield was obtained as one In fact, however, the quantum yield is less than one on account of the non-radiative transitions and abscrption events in which the total photon energy is converted into heat. In the present paper, the results of Refs 4 and 2 are used to calculate the quantum yield of the impurity fluorescence, taking the non-radiative transi tions into account. This is done by means of a perturbation operator V It is assumed that the zone of normal oscillations of the crystal may be divided into two classes: "optical" escillations which are in strong interaction with the localized electrons, and "acoustic" escillations which interact with them weakly Four states are considered: 1) the initial state; 2) the virtual state (electrons are in an excited state); 3) the final state (transition from the virtual state to the initial state with emission of radiation); 4) the final state (transition from the virtual state to the initial state without radiation). The formulas obtained by means of the terms of Ref. 2 show that the spectral relations derived in the earlier papers continue to hold in the present case also; only the quantum yield η is less than one. The temperature dependence of η is found to be given by $\eta = 1/(1+BT)$, which agrees qualitatively with experimental Card 2/3

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Consideration of Non-radiative Transitions S/181/60/002/008/051/052/XX in the Theory of Impurity Fluorescence B006/B070

observations. A. Ye. Marinchuk is mentioned There are 10 references. 6 Soviet, 2 German, 1 US, and 1 British

ASSOCIATION: Kishinevskiy gosudarstvennyy universitet (Kishinev State University)

SUBMITTED: January 19. 1960

Card 3/3

143600

2209, 1158, 1144

S/051/60/009/005/006/019

E201/E191

AUTHORS:

Perlin, Yu.Ye., and Palistrant, M.Ye.

TITLE:

On the Theory of Photoionization of F-Centres

PERIODICAL: Optika i spektroskopiya, 1960, Vol.9, No.5, pp 608-614

Continuous states of the F-centre spectrum are interpreted by the present authors as polarons moving in the The probability of screened Coulomb field of an amion vacancy. photoionization by transitions from the F-centre ground state is calculated. It is shown that due to participation of phonons in phototransitions, the dependence of the ionization probability on the frequency of absorbed light is an asymmetrical curve with a The long-wavelength end of this curve is similar to the Pekar curve (Ref. 5) for phototransitions in the discrete spectrum; the short-wavelength end of the authors curve fall: more smoothly than the Pekar curve. With increase of temperature the magnitude of the ionization probability maximum decreases and the half-width of the probability curve increases.

Card 1/2

S/051/60/009/005/006/019 B201/E191

On the Theory of Photoionization of F-centres

The calculated frequency dependences of the ionization probability for a KCl crystal at 76 °K (curve 1) and 794 °K (ourve 2) are given in a figure on page 614. Curve l'is very similar to an experimental dependence reported by Inchauspe

There are 1 figure and 5 references: 4 Soviet and 1 English.

SUBMITTED: February 26, 1960

Card 1/2

PERLIN, Yu. Ye., Doc Phys-Math Sci -- "Quantum transitions in local centers of crystals." Len, 1961. (Acad Sci USSR., Phys-Tech U im A. F. Ioffe) (KL, 8-61, 225)

- 3 -

PERLIN, Yu.Ye.; KOVARSKIY, V.A.

Theory of impurity-sensitived continued and the sensitive sensi

Theory of impurity-sensitized scattering of slow polarons. Fiz.tver. tels 3 no.4:1031-1034 Ap :61. (MIRA 14:4)

1. Kishinevskiy gosudarstvennyy universitet i Kishinevskiy sel'skokhozyaystvennyy institut. (Electrons—Scattering) (Color centers)

S/181/61/003/008/023/034 B109/B202

24.3500

Perlin, Yu. Ye., Marinchuk, A. Ye., Kon, L. Z.

TITLE:

AUTHORS:

Theory of the thermoluminescence of impurity crystals

PERIODICAL: Fiz_ka tverdogo tela, v. 3, no. 8, 1961, 2401-2412

TEXT: The thermoluminescence which occurs upon the radiationless transition of an electron from a metastable level to an excited level of a luminescence center is studied by the perturbation theory of Wigner-Weisskopf. The authors attempted to explain the temperature and frequency dependences of thermoluminescence intensity. As an example, the authors discuss the decoloration of the F' band which is due to a tunnel effect and is accompanied by a luminescence of the F-band. The calculation is made with the aid of an adiabatic model of a localized electron. The lattice spectrum is assumed to consist of two branches: optical vibrations whose interactions with the electron are calculated in zero-th approximation, and acoustic vibrations which interact only weakly with the electrons. This interaction is regarded as perturbation and causes radiationless electron transitions. If thermoluminescence is regarded as a second-order quantum transition and if Card 1/6

?**7293** 8/181/61/003/008/023/034 Theory of the thermoluminescence ... B109/B202 the excited electron-vibrational states are considered as virtual states, the theory of Wigner-Weisskopf generalized by Yu. Ye. Perlin (FTT, II, 1915, 1960 and FTT, II, 1928, 1960) can be used; the solution of the time-dependent Schrödinger equation is formulated as superpositions a) of the wave function Ψ of the initial state which is described by the wave function ψ_3 of the metastable level and the occupation numbers $n_{\mathbf{x}}^{o}$, $N_{\mathbf{f}}^{o}$, b) of the wave function $\Psi_{\mathbf{v}}$ of the virtual states with the electron wave function ψ_2 of the unstable level and the occupation numbers n_x , N_f , c) of the wave function Ψ_1 of the final states with the electron wave function ϕ_1 of the ground state and the occupation numbers $n'_{\mathbf{x}}$, $N'_{\mathbf{f}}$. The corresponding probability amplitudes are co, cowhose values can be calculated from the Schrödinger equation. The probability $\mathbf{w}(\Omega)$ of the emission of a photon $h\Omega$ can be found by taking the statistical mean value of $|c_r|^2$ $t \rightarrow \infty$ in terms of the occupation numbers of the photon oscillators in the initial state. Using the results and the denotation of A. Ye. Marinchuk, Yu. Ye. Perlin (Izv. Mold. fil. AN SSSR, 3, (69), 57, 1960) the authors obtain

Card 2/6



Theory of the thermoluminescence ...

APPROVED FOR RELEASE: 06/15/2000

$$w(2) = \frac{|W_{rr}(0)|^3}{\hbar^4 \Gamma \omega^3} \exp\left(-\frac{\alpha_{23} + \alpha_{21}}{2} \operatorname{cth} \frac{\beta}{2}\right) \sum_{f} |A_f|^3 \tilde{N}_f^0 \int\limits_{-\infty}^{\infty} d\tau \times$$

$$\times \exp \left[i\tau \left(p_0 - \frac{\omega_f}{\omega}\right) + \frac{a_{33}}{2 \sinh \frac{\beta}{3}} \cos \left(\tau - \frac{i\beta}{2}\right)\right] \int_{-\infty}^{\infty} dt \times$$

$$\times \exp \left[itr + \frac{a_{11}}{2 \sinh \frac{\beta}{2}} \cos \left(t - \frac{t\beta}{2}\right)\right] \int_{0}^{\infty} dt' \times$$

$$\times \exp \left[-\frac{\gamma}{\omega} t' - \frac{2ib}{\sinh \frac{\beta}{2}} \sinh \left(\frac{i\tau + \beta}{2} \right) \sin \frac{t}{2} \cos t' \right]$$

(2.14)

(2.15),

$$\times \exp \left[-\frac{\gamma}{\omega} t' - \frac{2ib}{\sinh \frac{\beta}{2}} \sinh \left(\frac{i\tau + \beta}{2} \right) \sin \frac{t}{2} \cos t' \right],$$

where

$$a_{11} = \sum_{i} (q_{13} - q_{11})^{2}; \qquad b = \sum_{i} (q_{13} - q_{13})(q_{11} - q_{12})$$

Card 3/6

CIA-RDP86-00513R001240110006-6"

Theory of the thermoluminescence ...

S/181/61/003/008/023/034 B109/B202

$$r\omega = \Omega - \Omega_{21}$$
 (2.16),
 $a_{32} = \sum_{x} (q_{x3} - q_{x2})^2$ (1.13),

$$A_{f} = \sqrt{\frac{1}{2D_{1}/2\omega_{f}}} \int \psi_{2}(\mathbf{r}) V_{f}(\mathbf{r}) e^{i(\theta,\mathbf{r})} \psi_{2}(\mathbf{r}) d\mathbf{r}, \qquad (1.12),$$

D - prystal density, L^3 - crystal volume. If $\zeta(\Omega)$ is the spectral frequency density then

$$\int w(\Omega) \rho(\Omega) d\Omega = \sum_{r \ge -\frac{\Omega_m}{m}}^{\infty} w_r \qquad (2.19)$$

follows from (2.14), ... the emission spectrum consists of equidistant lines. For w the authors give the following expression:

$$w_{r} = \exp\left(-\frac{\alpha_{21}}{2} \operatorname{cth} \frac{\beta}{2}\right) \sum_{k=-\infty}^{\infty} (-1)^{k} I_{k}^{2}(\xi) I_{r+k} \left(\frac{\alpha_{21}}{2 \operatorname{sh} \frac{\beta}{2}}\right) e^{-(r+k)\frac{\beta}{2}}. \quad (2.32)$$

Card 4/6

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theory of the thermoluminescence ...

S/181/61/003/008/023/0 B109/B202

This formula describes the frequency dependence of the intensity of thermoluminescence. For T \rightarrow 0, (2.32) changes into the formula for the probability of the light emission of an excited center after phonon relaxation. This formula was obtained by S. I. Tekar (ZhETF, 22, 641, 1952). For T / C the spectrum of the emitted radiation is a superfessit on of the curves for T = 0. It is bell-shaped and has a maximum at r = -1. 1/2.

The given formulas are applied to the case of thermoluminescence which occurs in a colored alkali halide crystal during the reaction $F^t \to F^t$. Table 1 gives the quantities $|J_f|^2$, and b for KCl as a function of the

distance R between the two centers; $a_{21} = 44.6$, $p_0 = -7$. Table 2 gives the decay rate \(\ \) of the centers as a function of R. Hence, at lower temperature

thermal ionization is less important. Thus, only the tunnel effect may cause a decoloration of the F' band. With increasing temperature the c. nditions are changed: According to A. G. Cheban (Opt. i spektr., \underline{X} , 495, 1951) the probability of thermal ionization at $T=300^{\circ} K$ is already ap, roximately 109 sec-1 and is thus of the same order of magnitude as the tunnel effect. There are 1 figure, 2 tables, and 11 references: | Seviet

Card 5/6

27293

Theory of the thermoluminescence ...

5,181,761/603/008/623/034 B109/B201

and 2 non-Soviet.

ASSOCIATION: Kishinevskiy gosudarstvennyy universitet Institut fizik: i matematiki Mold. fil. AM SSSR (Kishinevskiy State University

Institute of Physics and Mathematics AS USSR)

SUBMITTED:

March 16, 1961

R. Å	(<i>Īf</i>)*	a 30	1.61
5	0.48	30.6	7.53
7	0.21	43.2	9.05
10	0.062	57.1	8.13
15	0.32 · 10 ⁻²	70	7.2
20	0.37 · 10 ⁻⁴	73.1	7.0
90	0.59 · 10 ⁻⁸	80	6.9

R. A	Γ̄, cex1 (T = 30° K)	$\overline{\Gamma}_i \text{ cen.}^{-1}$ $(T = 300^{\circ} \text{ K})$	
5	2.3 · 10 ¹²	2.0 · 1013	
7	2.0 · 10 ¹⁰	1.8 · 1011	
10	4.3 · 10 ⁷	0.96 · 1011	
15	5 · 10 ⁴	0.89 · 109	
20	49 · 10	0.78 · 107	
30	1.0 · 10 ⁻³	0.65 · 169	

Table 1 Card 6/6

Table 2

S/051/60/008/03/018/038 E201/E191

AUTHOR:

Perlin, Yu.Ye.

Un the Problem of the Energy Spectrum of an F-Centre in

TITLE:

the Continuum Theory

PERIODICAL: Optika i spektroskopiya, 1960, Vol 8, Nr 3, pp 386-393 (USSR)

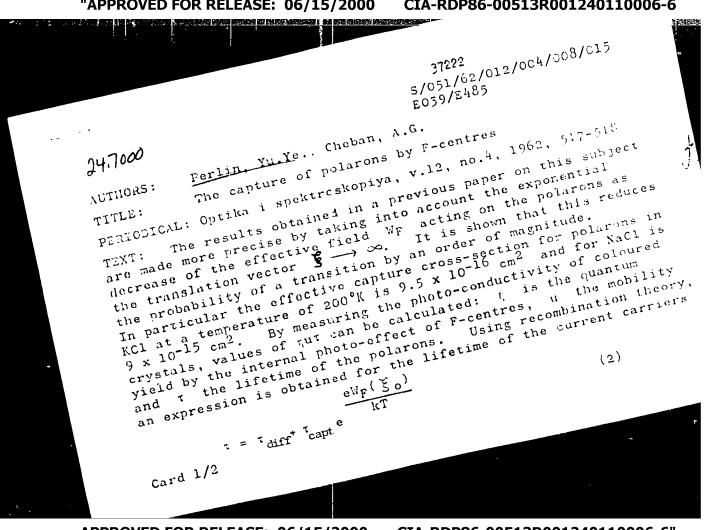
ABSTRACT: Within the framework of the continuum theory of an F-centre the author compares two types of variational functions. It is shown that functions of the first type (hydrogen-like model) give better results in calculation of the ground-state energy. For all the excited states functions of the second type (polaron model) yield more exact values of the energies. This indicates that in the theory of the excited states of an F-centre and other impurity centres in ionic crystals it is necessary to allow for the collective motion of phonons caused by the

Card 1/1

"excess" electrons (polaron effect). The paper is entirely theoretical. There are 7 references, of which 6 are Soviet and 1 translation from English.

SUBMITTED: July 8, 1959

CIA-RDP86-00513R001240110006-6 "APPROVED FOR RELEASE: 06/15/2000



\$/051/62/012/004/008/015 E039/E485

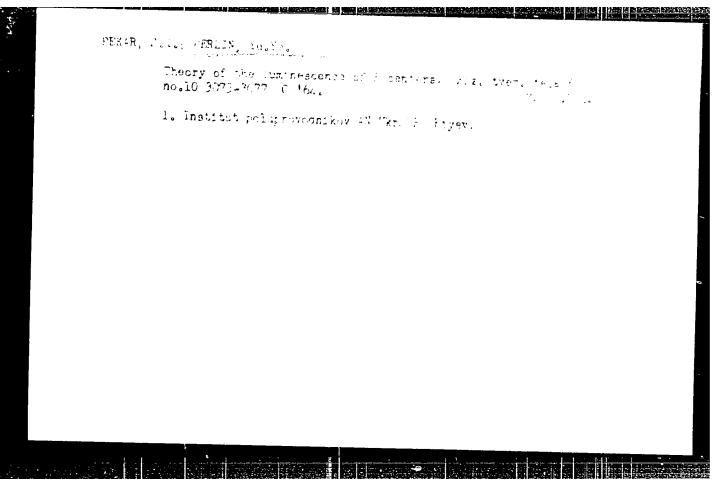
The capture of polarons ...

where $\tau_{\rm diff}$ - diffusion time of polarons to F-centres; $\tau_{\rm capt}$ - ligetime of a carrier with respect to quantum transitions at discrete levels; $W_F(\)$ - potential at "capture point".

At $-\frac{\sqrt{\sigma_\pi}}{10}$, this value is practically equal to zero in the low temperature region investigated where $\eta=1$. Values of ut are tabulated where τ is calculated from Eq.(2) and u from a formula by Pekar. These values of ut show good agreement with well-known experimental data for crystals of KCl and NaCl. There is 1 table.

SUBMITTED: May 18, 1961

Card 2/2



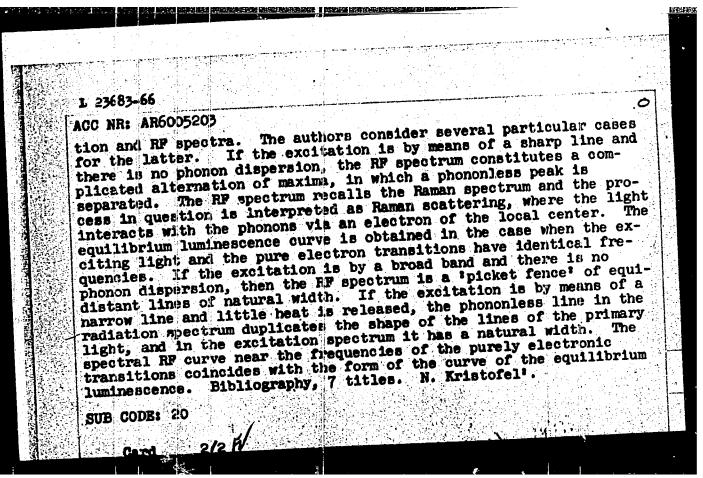
PHRLIN, Yu.Ye.; CHEBAN, A.G.

The K-absorption band in ionic crystals. Uch.zau.Kish.um. 69:1-2 *64. (MIRA 18:12)

1 10730 CZ mind 1 / 6 mm/ 1 mm/ 1	
1 12788-66 EWT(1)/T IJP(c) WM/GG	
ACC NR: AP5026618 SOURCE CODE: UR/0056/65/049/004/1237/1247	
AUTHORS: Gifeysman, Sh. N.; Perlin, Yu. Ye.	200
ORG: Kishinev State University (Kishinavality	
universitet	
TITLE: Impurity light absorption for arbitrary electron phonon	
SOURCE: Zhurnal eksperimental noy i teoreticheskoy fiziki, v. 49, no. 4, 1965, 1237-1247	
TOPIC TAGS: electron interaction, light absorption, phonon interaction, impurity center, thermodynamic function, Coulomb interaction, ionic crystal	
ABSTRACT: The method of arbitrary coupling, previously developed by one of the authors (Perlin, with I. Ya. Ogurtsov, FTT v. 7, 1467, 1965) for the calculation of thermodynamic functions of the impurity-phonon system, is extended in the present article to include the theory of impurity absorption of light. In this method, the spectral admittance of the	
by continual integration along the trajectories. In the zeroth approximation the system is described by a trial Lagrangian in which both the	
[Cont 1/2	

12788-65 ACC NR. AP5026618 Coulomb and the electron-phonon interactions are imitated by elastic forces. The need for this approach is dictated by the fact that the adiabatic approximation becomes inaccurate when the Bohr frequency of the pure electronic transition and the frequencies of the impurity-active phonons have the same order of magnitude, as is the case for shallow traps of the P-center type in some ionic crystals. The intensity of the no-phonon and one-phonon impurity absorption peaks is derived in first approximation for such traps. The formulas obtained are valid in the case of small heat release and arbitrary electronphonon coupling. The temperature shift of the impurity absorption lines is also calculated. Orig. art. has: 1 figure, 51 formulas, and SUEM DATE: 30Apr65/ 20/ SUB CODE: NR REF SOV: 009/ OTH REF: 007

IJP(c) ENT(I) SOURCE CODE: UR/0058/65/000/009/D062/D063 ACC NR: AR6005203 Rozenfel'd, Yu. B. Perlin, Yu. Ye. AUTHORS: On the theory of resonance fluorescence of local centers ${\cal B}$ TITLE: SOURCE: Ref. zh. Fizika, Abs. 9D492 Uch. zap. Kishinevsk. un-t, no. 75, 1964, 1-11 REF. SOURCE: TOPIC TAGS: fluorescence spectrum, phonon equilibrium, luminescence center, resonance line, excited electron state, line shape, electron transition TRANSLATION: 2/ The authors develop a general theory of impurity resonance luminescence in the case wher there is no time for phonon equilibrium to become established in the excited electron state. A modernized Wigner-Weisskopf method is used. Account is taken of the displacements of the equilibrium positions of the phonon coordinates. The change in the elastic constants is disregarded. Formulas are obtained for the probabilities of the resonance fluorescence (RF) and are used to obtain expressions describing the form of the excita-Card



L 04753-67 EWT(1)/EWT(m)/EWP(e)/EEC(k)-2/T/EWP(k) IJP(c) WG/WH

ACC NR. AP6025948 SOURCE CODE: UR/0051/66/021/001/0013/0018

AUTHOR: Tsukerblat, B. S.; Perlin, Yu. Ye.

ORG: none

TITLE: On theory of multiphonon nonradiating transitions between the local states of dissimilar multiplicity. II. Quantum yield of ruby luminescence on the R-line (low temperatures)

SOURCE: Optika i spektroskopiya, v. 21, no. 1, 1966, 13-18

TOPIC TAGS: electron transition, nonradiative transition, transition probability, transition radiation, ruby laser, ruby optic material, phonon, phonon interaction, crystal lattice energy, crystal lattice parameter, crystal lattice vibration

ABSTRACT: The authors use the adiabatic method to calculate the probability of a multiphonon nonradiation transition $^{2}E_{g}$ + $^{4}A_{2g}$ and the quantum yield of ruby luminescence of the R-line. The interaction with optical and acoustical vibrations of the crystal lattice is taken into account. The investigated transition corresponds to the operating levels of ruby lasers and occurs in the second order of perturbation theory via the virtual state $^{4}T_{2g}$. Other virtual states can be neglected because of their low probability. Taking in account only the strongest interaction of $^{2}E_{g}$ and $^{4}A_{2g}$

Card 1/2

UDC: 535.37 : 553.824.01

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ACC NR: AP6025948

electron energy levels with the wholly symmetrical lattice vibrations in the crystal, the adiabatic electron-oscillatory function is given by

$$\Psi\left(\Gamma\right) \stackrel{i}{\leftarrow} \psi_{\Gamma}\left(\mathbf{r}\right) \prod_{\mathbf{k}\sigma} \Phi_{N_{\mathbf{k}\sigma}}\left(q_{\mathbf{k}\sigma} - q_{\mathbf{k}\uparrow}^{i}\uparrow\right) \prod_{\mathbf{n},\sigma} \Phi_{n_{\mathbf{k}\sigma}}\left(q_{\mathbf{n}\sigma} - q_{\mathbf{n}\uparrow}^{i}\uparrow\right),$$

where $\psi_{\Gamma}(r)$ is the electron wave function from the crystal field theory; ϕ_n are wave functions of the harmonic oscillators; indices k, s identify wave vectors and zones of the acoustical phonons; z, σ are correspondingly related to the photons. The $q_{k\Gamma}^{(r)}(\lambda=k, z, \rho=s, \sigma)$

represent the displacements of the equilibrium for the oscillators in the phonon field due to electron-phonon interactions. The authors proceed from the standard expression for the probability of multiphonon nonradiation transitions, derived by the authors in an earlier work, to develop an expression which accounts for both the phonon and photon interactions in the crystal. Based on this expression the probability of nonradiation transition ${}^2E_g + {}^4A_{2g}$ at low temperatures has a value of 1.4 sec 1. Considering the lifetime of the E_g state to be 4.3·10⁻³ sec the quantum yield, in agreement with the experimental data, is practically equal to unity. Orig. art. has:

SUB CODE: 20/

SUBM DATE: 14Dec64/

ORIG REF: 004/

OTH REF: 002

Card 2/2 18

ACC NR: A.P6036979

(A, N)

SOURCE CODE:

UR/0181/66/008/011/3324/3334

AUTHOR: Perlin, Yu. Ye.; Gifeysman, Sh. N.

ORG: Kishinev State University (Kishinevskiy gosudarstvennyy universitet)

TITLE: Diamagnetism of bound polarons

SOURCE: Fizika tverdogo tela, v. 8, no. 11, 1966, 3324-3334

TOPIC TAGS: polaron, diamagnetism, impurity center, ionic crystal, electron spin, magnetic susceptibility, ionization

ABSTRACT: This is a continuation of earlier work (FTT v. 7, 1467, 1965 and elsewhere) dealing with hydrogenlike local centers in ionic crystals. The present article extends the earlier results to include the case of a weak homogeneous external magnetic field and the case of a bound polaron. Inasmuch as the Lagrangian formalism of quantum mechanics has been developed only for systems having a classical Lagrangian, the present investigation is limited to diamagnetic effects of the localized electron which are not coupled to the spin. The partition function of an impurity semiconductor in a magnetic field is represented in the form of a Feynman functional integral over the trajectories of the localized electron. In the zeroth approximation, the system is described by a trial Lagrangian in which the Coulomb terms are imitated by elastic interactions. The dependence of the trial-Lagrangian parameters on the magnetic field is disregarded. In the first approximation in the difference between the true and trial action the authors calculate the free-energy correction which is

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ACC NR. AP6036979

quadratic in the magnetic field intensity. Formulas for the diamagnetic susceptibility of the local centers is the function of the crystal parameters and the temperature are obtained for the limiting cases of strong and weak local coupling. In both limiting cases, allowance for the polaron effect leads to a decrease in the diamagnetic susceptibility compared with the hydrogenlike susceptibility by a factor 1.5 - 2 times. A formula is also obtained for the dependence of the ionization energy of the local center on the magnetic field intensity. Orig. art. has: 43 formulas and/1 table.

SUB CODE: 20/ SUBM DATE; 07Apr66/ ORIG REF: 008/ OTH REF: 013

Card 2/2

ACC NR AR7000872

SOURCE CODE: UR/0058/66/000/009/E074/E074

AUTHOR: Perlin, Yu. Ye.; Kharchenko, L. S.

TITLE: Quantum yield of the thermo-optical ionization of local centers

SOURCE: Ref. zh. Fizika, Abs. 9E605

REF SOURCE: Uch. zap. Kishinevsk. un-t, v. 80, 1965, 51-56

TOPIC TAGS: quantum yield, thermal optic effect, ionization, phonon

ABSTRACT: The quantum yield (η) of thermo-optical ionization of electrons from deep local levels into the conduction band was calculated, on the assumption that the non-radiative transfer of electrons from excited local levels to the conduction band occurs faster than the return of the phonon subsystem to a state of equilibrium (nonequilibrium thermo-optical process). It is shown that with sufficiently high Stokes losses, when radiative transfers from the nonequilibrium intermediate state to the basic state can be neglected, the probability of nonequilibrium thermooptical transfer is greater than the probability of an equilibrium transfer. In

Cord 1/2

this process the dependence of quantum yield on temperature is basic described by the $\exp(-\Delta E/kT)$, where the activation energy of mainly a function of light frequency. A calculation was made for a casimple zone, and a local center is examined in an adiabatic approximation of configurational curves is used to determine the electron-phrelationship. V. Tulvinskiy. [Translation of abstract]	f AE is
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ACC NR: AP7005333

SOURCE CODE: UN/ONS1/66/00 1/012/2-2-15-33

AUTHOR: Perlin, Yu. Ye.; Rozenfel'd, Yt. B.; Toukerblat, B. S.

ORG: Kishinev State University (Kishinevskiy gosudarstvennyy universitet)

TITLE: On the nature of the optical impurity absorption bands and luminescence of crystals activated with rare-earth ions

SOURCE: Fizika tverdogo tela, v. 6, no. 12, 1966, 3490-3499

TOPIC TAGS: activated crystal, luminescence, absorption band, light absorption, impurity center, Stark effect, phonon interaction, electron interaction

ABSTRACT: In view of the fact that the classical theory of the crystalline field, which takes into account the Stark splitting of the levels of the impurity ion by the quenched lattice, is insufficient for the interpretation of the spectra of impurity tabsorption and luminescence of trivalent ions of rare-earth elements (\mathbb{TR}^{3+}) in crystals of the MeF2 type, the authors employ the theory of impurity light absorption and luminescence in crystals, developed by M. A. Krivoglaz and S. I. Pekar (Trudy, Fmysics interaction on the shape of the impurity absorption or luminescence spectra in the the crystals CaS2, BeF2; and SrF2 activated with Ce3+. The interaction between the of the crystal is taken into account in the adiabatic approximation. It is shown that

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ACC NR: A17005833

the presence of local oscillations can lead, under certain conditions, to a periodic distribution of intensity in the optical spectrum. The parameters of the optical bands are calculated and a level scheme for the Ce³⁺ in the cubic crystalline field is presented. The theoretical calculations are compared with the experimental data of A. A. Kaplyanskiy et al. (Opt. i spektr. v. 14, 664, 1963) and reasons for some discrepancies are indicated. The authors thank S. I. Pekar, A. A. Kaplyanskiy, and B. Z. Malkin for useful discussion. Orig. art. has: 1 figure, 30 formulas, and 1

SUB CODE: 20/ SUBM DATE: 25Mar66/ ORIG REF: 011/ OTH REF: 004

Card 2/2

l 26765=66 BMI(1)/BMI(n)/IIJP(c) JD/JG/GG ACC NR. AT6005613 UR/2837/64/069/000/0001/0002 AUTHOR: Perlin, Yu. Ye.; Cheban, A.G. ORG: Kishinev State University, Kishinev (Kishinevskiy Gosuniversitet TITLE: On the problem of K-band absorption in ionic crystals SOURCE: Rishinev, Universitet, Uchenyye zapiski, v. 69, 1964, 1-2 TOPIC TAGS: crystal, ionic crystal, alkalai halide , crystal absorption , absorption band, polaron, ionization, K band, L band ABSTRACT: Previous work of the authors and others is reviewed to establish the theory that the M-absorption band can be related to photo-transitions of electrons in crystals from the ground F-center state into a polaron state. Certain computed and experimentally determined parameters (location of maxima of 12, and halfwidthm) of the F-photolonization curves invoking creation of polarons and K-bands are tabulated for the ionic alkali-halide crystals: NaCl, KCl, KBr, EI, RbCl, RbBr, RbI. Good correspondence the experimentally determined with the computed data is considered as supporting the hypothesis of phototransition into ionized polaron states as the origin of the Kband in alkali halide crystals. Interpretation of the L-bands requires additional tenes gy zone data evaluation for the alkali halide crystals. Orig. art. has: I table. יו SUBM DATE: ORIG REF: 003 20/ Home /

L 31493-66 ENT(1)/T IJP(c)

ACC NR: AP6013022

SOURCE CODE: UR/0051/66/020/004/0657/0660

经现金点

AUTHOR: Perlin, Yu. Ye.; Kovarskiy, V. A.; Tsukerblat, B. S.

50

ORG: none

 \mathcal{B}

TITLE: Contribution to the theory of many-phonon nonradiative transitions between local states of different multiplicity. I.

SOURCE: Optika i spektroskopiya, v. 20, no. 4, 1966, 657-660

TOPIC TAGS: nonradiative transition, spin orbit interaction, phonon interaction, electron interaction, SPINSISTEM

ABSTRACT: The authors analyze many-phonon nonradiative transitions between levels of different multiplicity within the framework of the adiabatic approximation. The <u>electron-phonon interaction</u> is assumed to be small compared with the spin-orbit interaction, and the non-adiabaticity operator is treated as the perturbation, using a method described by the authors elsewhere (FTT v. 4, 1936, 1962; Usp, fiz. nauk v. 80, 553, 1963). Possible mechanisms of many-phonon transitions are discussed, and a general formula is obtained for the probability of nonradiative transition with spin flip due to the spin-orbit interaction. The upper limiting case, when the spin-orbit interaction is small compared with the electron-

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VDC: 535.330: 548.0

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phonon interaction is considered in detail. The matrix element for the spin-orbit interaction is then regarded as the perturbation. Different expressions for the nonradiative transition are then obtained, depending on whether the electronic part of the matrix element of this transition vanishes or not. Orig. art. has: 17 formulas.

SUB CODE: 20/ SURM DATE: 14Dec64/ ORIG REF: 004/ OTH REF: 002

Card 2/2 mc

PERLIN, Yu.Ye.; ROZENFEL'D, Yu.B.

Theory of resonance fluorescence of local centers. Uch. zap. Kish. un. 75:1-11. 164. (MIRA 18:10)

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L 5612-66 EHA(k)/FED/ENT(1)/ENP(e)/ENT(m)/EEC(k)-2/ENP(1)/T/ENP(k)/ENA L 5612-66 EHA(k)/FED/ENT(1)/ENP(e)/ENT(m)/EEC(k)-2/ENP(1)/T/ENP(k)/ENA L 5612-66 EHA(k)/FED/ENT(1)/ENP(e)/ENT(m)/EEC(k)-2/ENP(1)/T/ENP(k)/ENA ACC NR: AP50274(5 SCTE/IJP(c) WG/NH SOURCE CODE: UR/0181/65/007/011/32	(n)-2/EMA(h) 78/2288
ACC NR: AP50274(5 My 50 AUTHOR: Tsukerblat, B. S.; Perlin, Yu. Ye. 47 55	64 B
AUTHOR: Tsukerblat, B. S.; Italian and State University (Kishimevskiy gosudrastvennyy university ORG: Kishimev State University (Kishimevskiy gosudrastvennyy university ORG: TITLE: On the theory of nonradiative transitions involving several photography:	nons in
SOURCE: Fizika tverdogo tela, v. 7, no. 11, 1903, or position, ruby laser	25,44
TOPIC TAGS: phonon interaction, nonrelated to probability of nonradiative multi-	
sitions between the orbital tripletage of sitions with d^3 configuration) in an octahedral crystal field. This other ions with d^3 configuration) in an octahedral crystal field. This other ions with d^3 configuration process for population inversion is the "bottleneck" in the relaxation process for population inversion to the "bottleneck" in the relaxation process for population to the considered as a much faster rate). States of virious multiplicity are considered at a much faster rate). States of virious multiplicity are considered at a much faster rate). States of virious multiplicity are considered at a much faster rate). States of virious multiplicity are considered at a much faster rate of the Cr3 ion is small in comparison with the effect of the cr3 ion is small in comparison with the effect of the cr3 ion is small in comparison with the effect of the cr3 ion is small in comparison with the effect of the cr3 ion is small in comparison with the effect of the cr3 ion is small in comparison with the effect of the cr3 ion is small in comparison with the effect of the cr3 ion is small in comparison with the effect of the cr3 ion is small in comparison with the effect of the cr3 ion is small in comparison with the effect of the cr3 ion is small in comparison with the effect of the cr3 ion is small in comparison.	The spin-
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tion. The formula for the transition is derived in the adiabatic approximation, taking the Jahn-Teller effect into account. Since the introduction of a chromium ion in place of aluminum in the corundum lattice (Al203) does not generate quasimolecular (localized) vibrations, a small number of variables will not suffice to describe the electron-vibration system. The final results are considerably dependent on the parameters of the crystal. The greatest contribution to the probability of the transition is from processes with simultaneous participation of longitudinal and transverse acoustic phonous. The authors thank I. B. Bersuker and B. G. Vekhter for useful consultation. ()rig. art. has: 40 formulas, 1 table:

SUB CODE: SS/

SUBN DATE: 28hay65/

ORIG REF: 007/

OTH REF: 011

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EWT(1)/7 L 00767-66 IJP(c) GG UR/0181/65/007/005/1467/1474 ACCESSION NR: AP5012559 Ye.; Ogurtsov, I. Ya. TITIE: Vibrational method in the thermodynamics of impurity-phonon systems SOURCE: Fizika tverdogo tela, v. 7, no. 5, 1965, 1467-1474 TOPIC TAGS: partition function, crystal impurity, variational method, ionic crystal, P center, electron interaction, phonon interaction ABSTRACT: An expression is derived for the partition function of an impurity crystal, in the form of a continual integral over the trajectories of the localized electron. The approach is based on a procedure first employed by R. P. Feynman (Phys. Rev. v. 97, 660, 1955), consisting of replacing the exact Hamiltonian of the system by a trial Hamiltonian containing variational parameters. In this article, a trial action is introduced, which makes it possible to reduce the calculation of the thermodynamic functions of an ionic crystal with an F-center to the problem of finding the maximum of a function of three parameters. The problem is solved in the limiting case of strong coupling between the localized electron and the lattice defect, for an arbitrary electron-phonon coupling. Formulas are obtained for the internal energy of the system, for the energy of the ground state, and for the impurity specific heat. The results are compared with those obtained by a different Cord 1/2

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	ABSOCIATION: SUPPLITIED: (MR KEF SOV:	08Dec64	Elich.	00 00 : 005	t (<u>Kishinev State Uni</u> SUB CODE: SS,	GP
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ENT(1)/T/EEC(6)-2 [[J])(c)/ESD(gs)/ASD(g)-5/AS(mp)-2 ACCESS (ON NR) AP 4045726 0/0030/64/006/003/0615/0625 AUTHOR Pekar 8 I ; Perlin, Yu. Ye TITLE: Local electron centers in ionic orystals 2 SOURCE, Physical status solidi, v. 6, no. 3, 1964, 615-625 TOPIC TAGS: F center, energy level, is state, 2p state, P luminescence, level electron center, emission probability, absorption probability ABSTRACT: This papel analyzes the statements by W. Beall Fowler and D. L. Dexter (Phys. stat. sol. 2, 821, 1962; 3, 1865, 1962) about the F-center theory. It is shown that, in spile of Stokes shift and difference in the Frank-Condon matrix elements of absorption and emission transitions, Einstein's relation for expressing the emission probability in terms of light absorption probability can be employed without limitations or approximations. Formulas are derived for the integral integral integral and experimental results, even in the case of alkali-halide crystals, when the electron orbit realus r does not exceed the lattice constant s (r <a), is explained. Formulas which do not depend on the ψ -function of the electrons and are Card 1/2

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common to all local ments fective mass method, calc Formulas of the continuou compared with those based dependence, the formulas alkali-halide clystals. cance in KCl arises from in the continuous P-center is higher than that of the ASSOCIATION: Institut Semiconductors, Academy of universitet (Kishiney State	on experimental data. However, the assumption made by Fowler double level is discussed theory the energy level of 2p state. Orig. art. has:	ever, owing to the strong r- r a factor of 2 to 3) for er and Dexter that F-liming
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EWT(1)/EEC(b)-2 ASD(a)-5/AS(mp)-2/AFWL/ESD(gs)/ESD(t)/IJP(c) ACCESSION BR: AP4046622 8/0181/64/006/010/3073/3077 AUTHORS; Pekar, S. I., Perlin, Va. Ye. Contribution to the theory of F-center luminescence TITLE: SOURCE, Fizika tverdogo tela, v. 6, no. 10, 1964, 3073-3077 TOPIC TAIS: luminescence, luminescence polarization, wave function, phototransition adiabatic approximation ABSTRACT The authors get around some of the mathematical difficulties involved in the adiabatic approximation by using an approach which is only approximately adiabatic, wherein the state of the electron and the corresponding energy are assumed to be slow functions of the normal coording/tes but these functions are expanded in powers of the deviations from the normal positions of the equilibrium positions of the normal coordinates. An electron radiative 2p - 1s phototransit on in an F center is considered, with account Card 1/3

L 16175-65 ACCESSION NR: AP4046622 of the interaction between the electron and the polarization wibrations of the littice and of the changes in the equilibrium positions of the long during the phototransition. The radii of the electron states are assumed to be sufficiently large, and the polarization of the crystal by the electron is calculated continually. For the nuclear configuration corresponding to the initial (2p) state of the electron, the authors calculate the energy levels and the wave functions $\psi_{1s}^{2p}(r)$ and $\psi_{2p}^{2p}(r)$ of the 2p and 1s states, and also the Frank-Condon matrix element of the radiative phototransition between them. Specific calculations for the KC1 crystal show that the Frank-Condon matrix element of the coordinate for the radiative transition can be 1.5--2 times larger than for the absorptive transition. This result is just the opposite of what was obtained by Beall Fowler and Dexter (Phys. Rev. v. 128, 2154, 1962). Orig. art. has: 17 formulas and 1 table. Card

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PERLIN, YU. YE.

Yu. Ye. Perlin, "Theory of the Impurity Centers."

report submitted for the Conference on Solid State Theory, held in Moscow,

December 2-12, 1963, sponsored by the Soviet Academy of Sciences.

PERLIN, Yu.Ye.

Modern methods in the theory of multiphonon processes. Usp. fiz.
nauk 80 no.4:553-595 Ag '63. (MIRA 1619)

PERLIN, Yu.Ye.; CHERAN, A.G.

On the theory of field ionization of local states.

Fig. twer. tela 4 no.11:3220-3227 N '62. (MIRA 15:12)

1. Kishinevskiy gosudarstvennyy universitet.

(Quantum theory) (Ionization)

L 10067-6; EWT(1)/BDS/EEC(b)-2--AFFTC/ASD/ESD-3--IJP(C)
ACCESSION IR: AR3000371 S/0058/63/000/004/E060/E060

SOURCH: Rich. Fizika, Abs. 4E404

58

AUTHOR: Perlin, Yu. Ye; CHeban, A. F.

TITIE On the theory of the decay of excited color centers in an electric field.

CITED SCIRCE: Tr. po fiz. polumovomikov. Kishinevsk. un-t, vyp. 1, 1962, 3-14

TOPIC TAGE: color center decay, conductivity of semiconductors

TRANSIATION: The probability of disintegration of an excited F-center by an external electric field is calculated. It is assumed that the intensity of this field E is much small than critical, at which the polarons disintegrate. The calculation is made within the framework of the continual model of the F-center in the adiabatic approximation. Use is made of the fact that in the zero approximation the excited state of the F-center can be interpreted as the motion of an undeformed polaron in a Coulomb field which is distorted at small distances.

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L 10067-63

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The quadratic Stark shift is disregarded. The solution of the equation for the polaron is sought by a variational method. A critical distance R sub 0 is introduced in such a way, that when the distance R of the polaron to the center is smiller thin R sub 0 it is possible to neglect its interaction with the external field, while when R is greater than R sub 0 it is possible to neglect the difference between the potential energy of the polaron and its Coulomb energy, and the solutions for these two regions join together at R sub 0. Calculation of the turnel current throught the barrier is carried out in the quasi-classical approximation. At absolute zero the quantum yield of the photo effect for K Cl is practically equal to zero up to E equals 2.6 times 10 sup 6 v/cm. Extrapolation to larger fields shows that in the narrow field interval from 2.6 times 10 sup (to 2.8 times 10 sup 6 v/cm the quantum yield increases to unity. The experimental curve is similar in shape to the theoretical curve, but the quantum yield approaches unity even it E equals 300,000 v/cm. The reason for the discrepancy may be related both with the errors in the calculation and the inhomogeneity of the external field, as well as with the sufficiently low temperature at which the experiments are performed. E. Nagayev

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PEKAR, S.I.; PERLIN, Yu.Ye.

Lifetime of excited F-centers. Zhur. eksp. i toer. fiz. 49 no.3:

1. Institut poluprovodnikov AN UkrSSSR.

(Ionic crystals) (Quantum theory)

1108-1110 162.

PERLIN, Yu.Ye.; CHEBAN, A.G.; TSUKERBLAT, B.S.

Theory of the capture and scattering of polarons by F-centers.

(MIRA 15:7)

Uch. zap. Kish. un. 49:11-18 '61. (MIRA 15:7)

(Electrons--Scattering) (Electrons--Capture) (Quantum theory)

l₁312l₁ **8/181/62/**004/011/024/049 B125/B186

AUTHORS:

Perlin, Yu. Ye., and Cheban, A. G.

TITLE:

On the theory of autoionization of local states

PERIODICAL:

Fizika tverdogo tela, v. 4, no. 11, 1962, 3220-3227

TEXT: The quasi-classical method of calculating the tunnel decay of a hydrogen-like atom (L. D. Landau and Ye. M. Lifshits. Kvantovaya mekhanika (Quantum mechanics), GITTL 1948) in a strong electromagnetic field is generalized to the case where the electron is localized in a Coulomb field distorted at small distances from the center. A polaron at rest and an excited F-center are taken as examples. When calculating the self-consistent state of a Pekar polaron (S. I. Pekar. Issledovaniya po elektronnoy teorii kristallov (Investigations on the electron theory of crystals), GITTL 1951) in a strong electric field it is necessary to consider the deformation of the self-consistent field besides the direct interaction between the electron and the external field. If text (for crystals of the alkali-halide type $\epsilon_{crit} \sim 3\cdot 10^6 \text{ v/cm}$), the deformation of the self-consistent field can be neglected. By this approximation the ionization Card 1/3

On the theory of autoionization ...

S/181/62/004/011/024/049 B125/B186

probability

$$w_{ion} = \alpha_0^3 \left(\frac{4}{\delta n^2}\right)^{2n\beta_0} f(\eta_0) \exp\left(-\frac{2}{3\delta n^3}\right), \qquad (1.25),$$

$$f(\eta_0) = \eta_0^{-2n_1} \exp\left[\eta_0\left(\frac{1}{n} - a_0\right)\right]. \tag{1.26}$$

can be derived from the electron wave function and the energy of the self-consistent state of the polaron at rest at 2 = 0, using the expression $\psi(\mathbf{r}) = \psi_1(\xi) \psi_2(\xi)$ for the axially symmetrical solution. Thereby $\alpha_0 = (5/16) \, \mu \, \mathrm{e}^2 \, \mathrm{c/h}^2$, $n = (-2E)^{-1/2}$, $n_2 = n\beta_2 = 1/2$, β_2 is one of the two constants for separation of variables. For η_0 of the parabolic coordinate $\eta = r-z$, the inequality $1 \le \eta_0 \le 2 |E|/E$ is valid. Thus, for KCl crystals the autoionization probabilities, given in \sec^{-1} , were 4.5, 7.1, 3·10⁵, 2.7·10⁷, 2.2·10⁶. These values correspond to $E = (0.6, 0.7, 0.8, 0.9, \text{ and } 9.1) \cdot 10^6 \, \text{v/cm}$. The critical field strength is apparently $\mathbf{f} \sim 8 \cdot 10^5 \, \text{v/cm}$. The autoionization probability for an excited F-center with a transition of a localized electron from the potential well into the free state is derived Card 2/3

On the theory of autoionization...

S/181/62/004/011/024/049 B125/B186

similarly by solving the wave equation for the electron wave function ψ_2 in the presence of an external homogeneous field. For ψ_2 an exponential expression with parabolic coordinates is used. The autoionization probabilities for an excited F-center were 6.2·10⁴; 9.5·10⁵; 5.0·10⁶; 1.7·10⁷ sec⁻¹ and 4.9·10⁹ for the **L**-values (2.2; 2.3; 2.4; 2.5; and 3.0)·10⁶ v/cm. The autoionization is considered to be the only possible reason for the decay of excited F-centers near absolute zero. There are 2 figures and 2 tables.

ASSOCIATION: Kishinevskiy cosudarstvennyy universitet (Kishinev State

University)

SUBMITTED:

June 25, 1962

Card 3/3

PERLIN, Yu.Ye.; CHEBAN, A.G.

Poleron capture by F-centers. Opt. 1 spektr. 7 no.4:517-518
Ap '62. (MRA 15:5)

(Color centers) (Quantum theory)

经各种股份的 医阿拉斯斯氏 医克里克斯氏 医阿里克斯氏炎 二十二次 计图像 医自己外向 1. 7. 1 s/056/62/043/003/055/065 B104/B102 14.3500 Pekar, S. I., Perlin, Yu. Ye. AU THORS: The lifetime of excited F centers TITLE: Zhurnal eksperimental noy i teoreticheskoy fiziki, v. . . , PURIODICAL: ne. 7(9), 1962, 1108-1110 TEXT: Calculated and experimental results for the radiation lifetimes of excited F centers are shown to be in good agreement, the former because based on the continuous theory of F centers of ionic crystals establishes by S. I. Pekar (Issledovaniya po elektronnoy teorii kristallov - Research into the electron theory of crystals, Gostekhizdat, 1951; ZhETF, 22, 541. 1952) and the latter obtained by R. K. Swank and F. C. Brown (Phys. Rev. Lett., 8, 10, 1962). The formula $\tau_{R}^{-1} = 4e^{2}n\Omega_{m}^{3}z_{21}^{2}/3\hbar c^{3}$ is deduced for the radiation lifetime. $\Omega_{\rm m}$ is the maximum frequency in the luminescence band, in is the refractive index. Results: Card 1/2

i -	The lifetime	of excited F centers	\$/056/62/043/003/05//06? B104/B102			أو عا		
· • •			K C l	KBr	KI.			
	10 ⁶ τ	(experimental), sec	0.59	1.1	2.1			
	10 ⁶ τ.	(calculated), sec	0.59	0.59	0.54			
	Thus theory justifies making not only qualitative but also quantitative statements as to the infrared luminescence bands and radiation lifetime of the F centers. ASSOCIATION: Institut poluprovodnikov Akademii nauk USSR (Institute in							
	SUBMITTED:	Semiconductors of the Academy of Sciences UkrSSR) : June 16, 1962						
	11 1	,						
	Card 2/2							

PERLINA, A. M.

"Water Softening in a City Water-Supply System." Sub 3 Jun 47, Moscow Inst of Engineers of Municipal Construction

Dissertations presented for degrees in science and engineering in Moscow in 1947.

SO: Sum.No. 457, 18 Apr 55

Land Tech See.

PERLINA, A. M.

Voshchenko, Z. S. and Perlina, A. M. "Soft waters in municipal economy and living conditions," San. Teknnika, Issue, 2, 1948, p. 19-44

SO: U-2888, Letopis Zmurnal'nykh Statey, No.1 , 1949

PERLINA, A.M.,
POPOV. I.P., kardidat biologichenkikh nauk; PERLINA, A.M., kandidat
tekhnicheskikh nauk; KASHAYRV, S.I.

Water moftening in Moscow laundries. Gor.khoz.Mosk. 28 no.8:22-23
Ag 154. (MIRA 7:9)

1. Starshiy tekhnolog Moskovskogo gorodskogo tresta prachechnykh (for Kushayev) (Mater-Softening) (Moscow-Laundries, Public) (Laundries, Public --Moscow)

BLIOKH, S.S., kandidet mediteinskikh nauk; PKRLINA, A.M., kandidet tekhnicheskikh nauk; KCZLOVA, N.L., inzhener

Bifectiveness of the new method of purifying drinking water (contract clarification). Gig. i sen. 22 no.1:70-72 Js '57. (MIRA 10:2)

1. In Nauchno-issledovatel'skogo senitarnogo institute imeni Brismans, Akademii kommunal'nogo khozyaystva imeni Pamfilova i Laboratorii Rublevskoy vodoprovodnoy stentsii.

(WATER SUPPLY, purification, contact clearing technic (Rus))

Prolina F/ 1/1

USSR Chemical Technology. Chemical Products

H-5

and Their Application

Water treatment. Sewage water.

Abs Jour: Referat Zhur - Khimiya, No 1, 1958, 1697

: Perlina A.M. Author

: Experience with the Use of Contact Clarifiers Title

Orig Pub: Sb.: Issledovaniya po vodopodgotovke. M., Gos. izd-vo po str-vu i arkhitekt., 1956, 86-89

Abstract: Results of tests of industrial contact clarifiers (CC) at a number of water supply stations. All the CC produce a lowering of the color and turbidity of the water to meet the GOST. In bacter-iological and hydrobiological indices the water is practically identical with water purified in conventional units. As concerns the zooplankton a somewhat inferior purification result has been

Card 1/2

USSR Chemical Technology. Chemical Traducts and Their Application
Water treatment. Sewage water.

Abs Jour: Referat Zhur - Khimiya, No 1, 1958, 1697

attained. Congulating agent dosages amount 20-55% of those used at conventional installations. With a turbidity of the initial water not above 4-5 mg/liter and a coloration not in excers of 28-30, the CC can operate without congulation. Minimum duration of operation cycle of CC is greater than the permissible limit in the case of rapid filters. The average yearly consumption of wash water is 3.5-9%; the maximal (during floods) is 10-18%. To control the horizontal shifting of gravel layers it is recommended to use distributing systems having a large resistance without horizontal compensation.

Card 2/2

PISEUROV, Pawel Ivanovich, prof., doktor tekhn.nauk; NAGORMOV, Mikolay
Ivanovich; FERLINA, A.M., red.; SHVEDOV, Yu.F., red.izd-va;
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PERLINA, A.M.; BALASHOVA, G.V.; GCRYAINOVA, G.S.

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SHUBERT, S.A.; PERLINA, A.M.; KULZHKHSKIY, V.I.; SIDENKO, T.K.; ALEKSANDROV, D.H.; SOKOLOV, V.F.; PAL'KOVSKAYA, L.N.; BRUK-LEVINSON, T.L.; BELYAKOVA, A.M.; KOZHEVHIKOVA, Ye.K.; AVRUSHCHENKO, R.A., red. izd-va; VOLKOV, S.V., tekhn.red.

[Weter purification for water supply to machine-tractor stations and state ferms] Ochisthe vody dlie vodosnebzhenite poselkov MTS i sovkhozov. Moskva, Izd-vo M-va kommun.khoz. RSFSR, 1957.

69 p. (MIRA 11:6)

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Paklilla, F.E., kandidat meditsinskikh nauk

Metastatic cerebral abscesses in children. Pediatriis no.3:74-77

Hetsitatic cerebral abacesses in children. Fediatrita do.): (447)
Hr 157. (MIRA 10:10)

1. Az 1-y geredskoy infektsionnoy bol'nitsy Dnepropetrovska (glavnyy vrach F.E.Byushko) i Detskoy klinicheskoy bol'nitsy iglavnyy vrach L.V.Volkova) (BRAIN--ABSCESS)

PERLINA, F.I., kandidat meditsinskikh nauk

Clinical peculiarities of acute poliomyelitis in adults. Sov.med. 21 no.5:81-84 My '57. (MIRA 10:7)

1. Iz kliniki infektsionnykh bolezney (zev. - prof. I.I.Levin)

Dnepropetrovskogo meditsinskogo instituta i 1-y gorodskoy infektsionnoy bol'nitsy (glavnyy vrach - F.K.Nyushko)

(POLIOMYELITIS

in adults, clin. aspects)

PERLINA, P.I., kand.med.nouk

Clinical aspects and diagnosis of aparalytic and abortive forms of polionyelitis. Sov.med. 21 no.12:41-46 D '57. (NIRA 11:3)

1. Is kafedry infektsionnykh bolezney (sav.-prof. I.I.Levin)
Dnepropetrovskogo meditsinskogo instituta i Pervoy gorodskoy
infektsionnoy bol'nitsy (glavnyy vrach F.K.Nyushko)
(POLIOHYELITIS
aparalytic & mild forms, diag. & clin, manifest. (Rus)

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Affections of the pervous system in antirebies vencination. Thursdiscondinger of industry and the pervous system in antirebies vencination. Thursdisconding the pervoy gorodsky infektslooney belinitary i Detakny klinicheskoy belinitary.

(DEMYOUS STADE, diseases, caused by relical various (Rus))

(RABIES, including to dis. (Rus))
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PERLIHA, F., I.

Etiology and clinical aspects of brain abscesses. Hov.khir.erkh. np.2:107-108 Mr-Ap 158 (MIRA 11:6)

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(ERA IN-ABSCESS)

IEVIE, I.I., prof. [decessed], PTRLINA, F.I. kand,med.nauk

Acute rheumatic meningitie. Vrech.delo no.6:573-577 Je '58

(NIBA 11:7)

1. Elinika infektionnyih bolesney (sav.- prof. I.I. Levin [decembed])

Inepropetrovskogo mediteinskogo instituta i Pervaya gorodskaya infektionnaya bol'nites.

(MENINGITIS)

(RHEUMATIC FEVEE)

PERLINA, F.I., kand. meditsinskikh nauk

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(NERVOUS SYSTEM--DISEASES)

PERLINA, P.I.

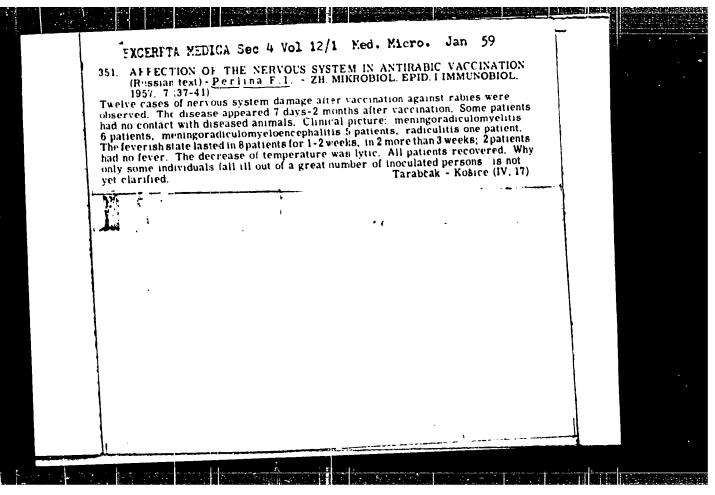
Role of the posterior radioles of the spinal cord in the regulation of trophic functions; osteourthropathy and changes in the gastro-intestinal tract in tabes dorsalis. Zhur. nevr.i psikh. 60 no.10: (MIRA 14:1) 1281-1290 '60.

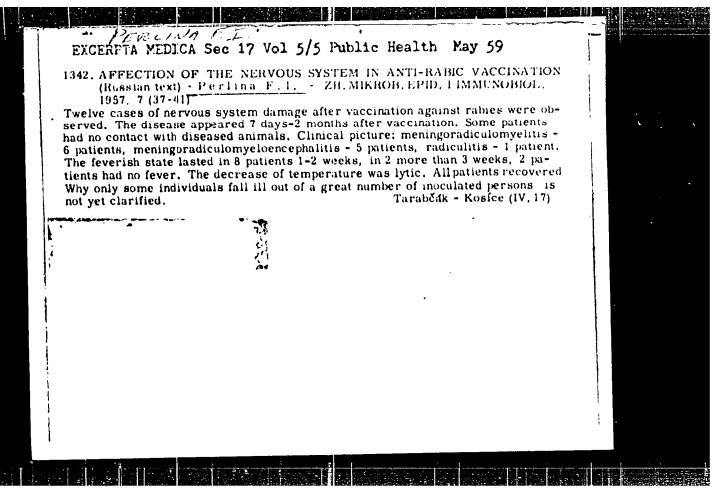
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(LOCOMOTOR ATAXIA) (NIGESTIVE ORGANS)

(BONES-DISEASES)

(JOINTS-DISEASES)





Acute paralysis of the facial nerve and policryclitis. Klinened.
36 no.9:119-124 S'58

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Dnepropetrovskogo meditsinskogo instituta i percoy gorodskoy
infektsionnoy bol'nitsy (glavnyy vrach F.K. Hyushko).

(POLIONYELITIS, compl.
acute facial nerve paralysis (Rus))

(FACIAL PARALYSIS, eticl & pathogen.
policryclitis (Rus))

PURLINA, F. M.

Perlina, F. M. "On the problem of distribunce to the cody system in peripheral trauma," Nevronatologiya i msikolatriya, 1949, No. 2, p. 32-3.

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	(Contd) certain parts of the brai	Repurts observation on cases with loss of reali- sation of a person's own body, his condition, the occurrence of a feeling of additional parts of the body, etc. Clinico-anatomical studies indi- cate existence of a close relationship of this	isturbances	Wesh/Medicine - Forve Diseases Medicine - Trauma - Complications and Sequels	FA 61/4975
61/13	o1/49790 Mar/Apr 19	reali- ton, the te of this this	of the Ferline,	Mar/Apr 49	

PERLINA L.P.

USSR/ Microbiology - Sanitary Microbiology.

F-3

Abs Jour: Ref Zhur-Biol., No 9, 1958, 38421

Perlina, L. P. Inst : Not given

Title: : Morphological and Biological Properties of Pyogenic Microorganisms Found in Maternity Homes.

Orig Pub: Sb. nauchn. rabot. Mold. otd. Vses. nauchn. o-va mikrobiol., epidemiol. i infektsionistov, 1956, No 1, 27-36.

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- 2. USSR (600)
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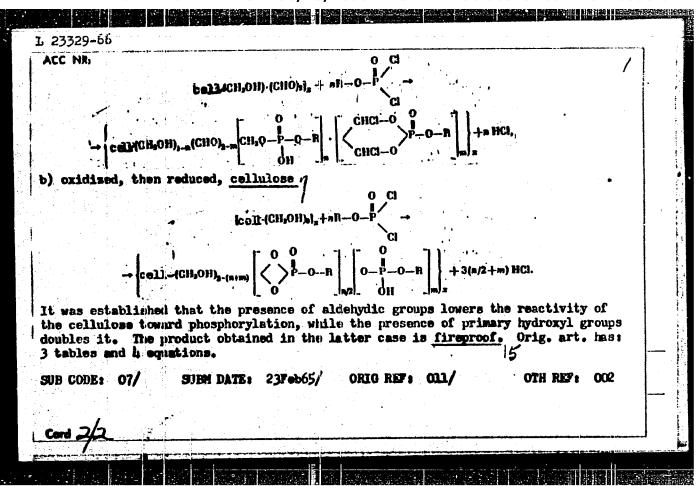
USMANOV, Kh.U.; PERLINA, R.V.

Determination of aldehyde and carboxyl groups in cellulose preparations. Uzb.khim.zhur. no.4:22-31 '61. (MIRA 14:8)

1. Institut khimii polimerov AN UzSSR. 2. Chlen-korrespondent AN UzSSR (for Usmanov).

(Cellulose) (Aldehydes) (Carboxyl group)

ACC NR. AP6005975 (A)	SOURCE CODE: UR/0190/66/008/002/0231/023h	
AUTHORS: Yuldashev, A.; Perlina, R.	V.; Sadykov, M. M.; Usmanov, Kh. U.	
ORG: Scientific Research Institute of (Nauchno-isaledovatel'skiy institut k	Chemistry and Technology of Cotton Cellulose himii 1 tekhnologii khlopkovoy tsellyulozy)	
TITIE: Phosphorylation of modified c chloroamydrides	ellulose preparations with phosphoric	
SOURCE: Vysokomolekulyarnyye soyedin TOPIC TAGS: <u>cellulose plastic</u> , phosp	eniya, v. 8, no. 2, 1966, 231-234 horylation, organic phosphorus compound	
present in position 2 and 3 of cellul with phosphoric dichloroanhydride (II ratio 1:11 = 1:3 were reacted in 30 m for 30 mimtes. The reactivity of th phosphorylation was determined from t Phosphorylation of the native cellulo by Wu. Mei-yen, T. A. Zharova, and Z.	nt of aldehydic or primary hydroxyl groups ose (I) upon the phosphorylation process of I) has been investigated. The reagents in the 1 of benzene and 10 ml of pyridine at 75-800 e various cellulose preparations toward he amount of P taken up during the reaction. se was described, and the mechanism was suggeste A. Rogovin (Zh. prikl. khimii, 35, 1820, 1962). lose proceeds according to the following schemes	
Cord 1/2	UDC: 661,728,87	

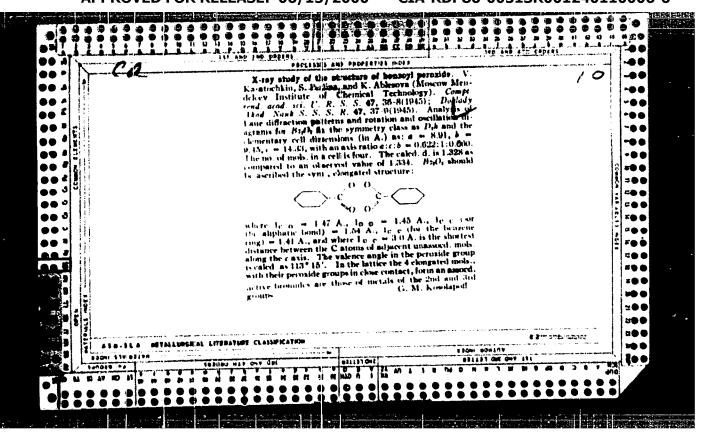


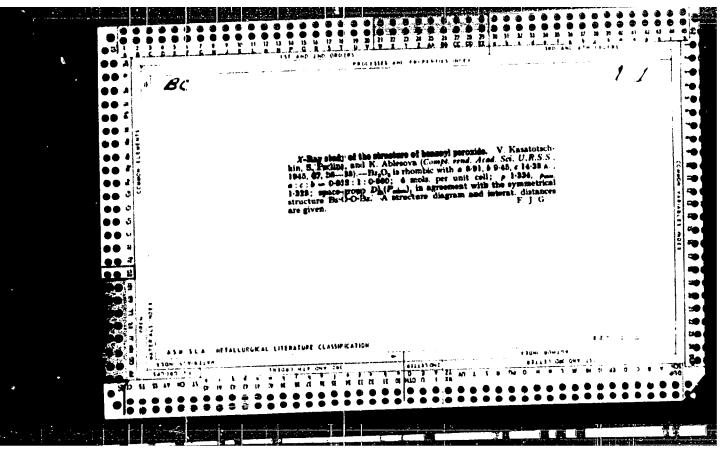
PERLINA, S.

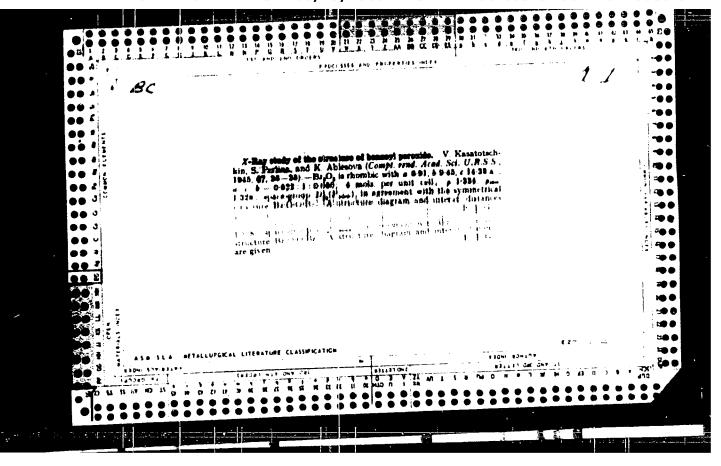
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OKUNEV, Yu.K., podpolkovnik, red.; MKDNIKOVA, A.N., tekhn.red.

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(DIARRHEA, in infant and child bacteriol, eticle in Poland)